Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100815\
Data File : BE090998.D
Acq On : 8 Oct 2015 11:57
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 09 03:17:49 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 07 17:11:43 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	167#	0.00
2		1,4-Dioxane	0.888	1.833	-106.4#	402#	-0.10
3		n-Nitrosodimethylamine	0.879	0.680	22.6	143	-0.13
4	S	2-Fluorophenol	1.045	1.136	-8.7	207#	-0.02
5	S	Phenol-d6	1.280	1.471	-14.9	215#	-0.02
6		bis(2-Chloroethyl)ether	1.524	3.741	-145.5#	425#	-0.05
7	I	Naphthalene-d8	1.000	1.000	0.0	172#	0.00
8	S	Nitrobenzene-d5	0.270	0.370	-37.0#	249#	-0.02
9		Nitrobenzene	0.289	0.430	-48.8#	286#	-0.01
10		Naphthalene	1.075	1.027	4.5	172#	-0.02
11		Hexachlorobutadiene	0.179	0.148	17.3	143	-0.02
12		2-Methylnaphthalene	0.634	0.687	-8.4	198#	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	176#	0.00
14	S	2,4,6-Tribromophenol	0.158	0.161	-1.9	240#	0.00
15	S	2-Fluorobiphenyl	1.355	1.323	2.4	185#	-0.01
16		Acenaphthylene	8.351	8.262	1.1	190#	0.00
17		Acenaphthene	1.334	1.266	5.1	178#	0.00
18		Fluorene	1.649	1.623	1.6	186#	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	185#	-0.01
20		4-Bromophenyl-phenylether	0.189	0.182	3.7	194#	-0.01
21		Hexachlorobenzene	0.970	0.198	79.6#	40#	-0.01
22		Pentachlorophenol	0.070	0.073	-4.3	302#	0.00
23		Phenanthrene	1.265	1.158	8.5	186#	-0.01
24		Anthracene	1.131	1.099	2.8	204#	-0.01
25		Fluoranthene	1.503	1.375	8.5	208#	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	195#	0.00
27		Benzidine	0.090	0.208	-131.1#	552#	-0.03
28		Pyrene	1.277	1.134	11.2	207#	0.00
29	S	Terphenyl-d14	0.631	0.625	1.0	221#	0.00
30		Benzo(a)anthracene	1.176	1.201	-2.1	221#	0.00
31		3,3'-Dichlorobenzidine	0.201	0.000	100.0#	0#	-20.84#
32		Chrysene	1.505	1.240	17.6	188#	0.00
33		Bis(2-ethylhexyl)phthalate	0.469	0.576	-22.8	375#	0.00
34		Indeno(1,2,3-cd)pyrene	1.212	1.292	-6.6	240#	0.00
35	I	Perylene-d12	1.000	1.000	0.0	229#	0.00
36		Benzo(b)fluoranthene	1.102	1.241	-12.6	288#	0.00
37		Benzo(k)fluoranthene	1.515	1.156	23.7	195#	0.00
38	C	Benzo(a)pyrene	1.022	1.112	-8.8	289#	0.00
39		Dibenzo(a,h)anthracene	1.009	1.212	-20.1	327#	0.00
40		Benzo(g,h,i)perylene	1.096	1.248	-13.9	301#	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 0